

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101819\
 Data File : BN008254.D
 Acq On : 18 Oct 2019 19:11
 Operator : JU
 Sample : K5178-09DL 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPF2DL

Manual Integrations
 APPROVED

mohammad
 10/22/2019 1:40:49 AM

Quant Time: Oct 19 00:05:04 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 23:47:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	3031	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	14409	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	9088	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	18444	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	17533	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	15882	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	361	0.01	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	1035	0.02	ng/ul	0.00
Target Compounds					Ovalue	
12) Phenanthrene	17.02	178	2876	0.053	ng/ul#	93
15) Fluoranthene	19.05	202	5406	0.076	ng/ul	94
17) Pyrene	19.41	202	5514	0.079	ng/ul#	94
18) Benzo(a)anthracene	21.17	228	3139	0.049	ng/ul#	90
19) Chrysene	21.22	228	3064	0.039	ng/ul#	92
21) Benzo(b)fluoranthene	22.72	252	4710m	0.091	ng/ul	
22) Benzo(k)fluoranthene	22.75	252	1748m	0.030	ng/ul	
23) Benzo(a)pyrene	23.27	252	2647	0.049	ng/ul#	24
24) Indeno(1,2,3-cd)pyrene	25.52	276	1518	0.024	ng/ul#	93
26) Benzo(g,h,i)perylene	26.18	276	1614	0.027	ng/ul#	67

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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